## Claims

We claim:

1. A compound of the structure

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wherein Y, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>1</sup>, C(R<sup>2</sup>)(R<sup>3</sup>), NR<sup>5</sup>, CH, O and S;

q is an integer of from 3 to 10;

A is selected from the group consisting of O, S,  $C(R^{16})(R^{17})$  and  $NR^{6}$ ;

E is selected from the group consisting of CH<sub>2</sub>, O, S, and

NR<sup>7</sup>;

J is selected from the group consisting of O, S and NR<sup>8</sup>;

T is selected from the group consisting of C(O) and  $(CH_2)_b$  wherein b is an integer of from 0 to 3;

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M is selected from the group consisting of  $C(R^9)(R^{10})$  and  $(CH_2)_u$ , wherein u is an integer of from 0 to 3;

L is selected from the group consisting of O, NR<sup>11</sup>, S, and (CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

X is selected from the group consisting of CO<sub>2</sub>B, PO<sub>3</sub>H<sub>2</sub>,

SO<sub>3</sub>H, SO<sub>2</sub>NH<sub>2</sub>, SO<sub>2</sub>NHCOR<sup>12</sup>, OPO<sub>3</sub>H<sub>2</sub>, C(O)NHC(O)R<sup>13</sup>.

C(O)NHSO<sub>2</sub>R<sup>14</sup>, hydroxyl, tetrazolyl and hydrogen;

W is selected from the group consisting of C, CR<sup>15</sup> and N; and

B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup>, R<sup>12</sup>, R<sup>13</sup>, R<sup>14</sup>, R<sup>15</sup>, R<sup>16</sup> and R<sup>17</sup> at each occurrence are independently selected from the group consisting of

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hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy,  $-N(C_1-C_3 \text{ alkyl})-C(O)(C_1-C_3 \text{ alkyl})$ ,  $-NHC(O)N(C_1-C_3 \text{ alkyl})$ alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub>alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub>alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>- $C_3$ )amino,  $-C(O)O-(C_1-C_3)$ alkyl,  $-C(O)NH-(C_1-C_3)$ alkyl,  $-C(O)N(C_1-C_3)$ alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup>, R<sup>12</sup>, R<sup>13</sup> R<sup>14</sup>, R<sup>15</sup>, R<sup>16</sup> and R<sup>17</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring: and wherein when M is C(R<sup>9</sup>)(R<sup>10</sup>), R<sup>9</sup> and R<sup>10</sup> taken together may form a ring; and wherein when A is NR<sup>6</sup> and at least one Y is CR<sup>1</sup>, R<sup>1</sup> and R<sup>6</sup> taken together may form a ring; or a pharmaceutically acceptable salt thereof;

with the proviso that when A is  $C(R^{16})(R^{17})$ , E is not  $NR^{7}$ .

25 2. A compound of claim 1 wherein

> A is  $NR^6$ ; E is NR<sup>7</sup>:

J is O;

M is  $C(R^9)(R^{10})$ ;

30 q is 4 or 5;

T is  $(CH_2)_b$  wherein b is 0;

L is  $(CH_2)_n$  wherein n is 0;

X is CO<sub>2</sub>B;

W is C or CR<sup>15</sup>;

R<sup>4</sup> is selected from the group consisting of aryl, alkylaryl, aralkyl, heterocyclyl, alkylheterocyclyl and heterocyclylalkyl; and R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup> and R<sup>15</sup> are independently selected from the group consisting of hydrogen and lower alkyl.

- 3. A compound of claim 1 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.
  - 4. A compound of the structure

wherein Y, at each occurrence, is independently selected from the group consisting of C(O), N,  $CR^1$ ,  $C(R^2)(R^3)$ ,  $NR^5$ , CH, O and S;

q is an integer of from 3 to 7;

T is selected from the group consisting of C(O) and  $(CH_2)_b$  wherein b is an integer of 0 to 3;

L is selected from the group consisting of O,  $NR^{11}$ , S, and  $(CH_2)_n$  wherein n is an integer of 0 or 1;

W is selected from the group consisting of C, CR<sup>15</sup> and N; and

B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup> and R<sup>15</sup> are independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl,

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heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl),
                              -NHC(O)N(C_1-C_3 \text{ alkyl})C(O)NH(C_1-C_3 \text{ alkyl}), -NHC(O)NH(C_1-C_6 \text{ alkyl}),
                             -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino.
                             alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl,
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                              -C(O)NH-(C_1-C_3)alkyl, -C(O)N(C_1-C_3)alkyl)_2, -CH=NOH, -PO_3H_2,
                             -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide.
                             cycloalkyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl,
                             aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl,
                             aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-
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                             C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and
                             -C(O)NH(benzyl) groups;
                            wherein B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup> and R<sup>15</sup> are
                                       unsubstituted or substituted with at least one electron donating or
                                      electron withdrawing group;
                            wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring;
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                                      and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring:
                                      and wherein when at least one Y is CR1, R1 and R6 taken together
                                                may form a ring;
                   or a pharmaceutically acceptable salt thereof.
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        5.
                  A compound of claim 4 wherein
                  q is 4 or 5;
                  W is C or CR<sup>15</sup>;
                  T is (CH_2)_b wherein b is 0;
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                  L is (CH_2)_n wherein n is 0;
                  R<sup>4</sup> is selected from the group consisting of aryl, alkylaryl, aralkyl,
                            heterocyclyl, alkylheterocyclyl and heterocyclylalkyl; and
                  R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup> and R<sup>15</sup> are independently selected from the
                            group consisting of hydrogen and lower alkyl.
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- 6. A compound of claim 4 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.
- 5 7. A compound of the structure

- wherein Y, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>1</sup>, C(R<sup>2</sup>)(R<sup>3</sup>), NR<sup>5</sup>, CH, O and S;
- q is an integer of from 2 to 5;
  - T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of 0 to 3;
  - L is selected from the group consisting of O,  $NR^{11}$ , S, and  $(CH_2)_n$  wherein n is an integer of 0 or 1;
- 15 R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>11</sup> and R<sup>18</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups; and
  - B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>9</sup> and R<sup>10</sup> are independently selected from the group consisting of hydrogen, halogen, halkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl,
- 25 -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl,

heterocycloyl, carboxy,  $-N(C_1-C_3 \text{ alkyl})-C(O)(C_1-C_3 \text{ alkyl})$ ,  $-NHC(O)N(C_1-C_3 \text{ alkyl})C(O)NH(C_1-C_3 \text{ alkyl}), -NHC(O)NH(C_1-C_6 \text{ alkyl}),$ -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl,  $-C(O)NH-(C_1-C_3)alkyl, -C(O)N(C_1-C_3alkyl)_2, -CH=NOH, -PO_3H_2,$ -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein B, R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup> and R<sup>18</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring: and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring; and wherein when at least one Y is CR<sup>1</sup>, R<sup>1</sup> and R<sup>6</sup> taken together may form a ring; or a pharmaceutically acceptable salt thereof.

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8. A compound of claim 7 wherein R<sup>18</sup> is selected from the group consisting of hydrogen, alkyl, aryl, aralkyl, cycloalkyl, alkylheterocyclyl, heterocyclylalkyl and heterocyclyl;

T is (CH<sub>2</sub>)<sub>b</sub> wherein b is 0;

- L is  $(CH_2)_n$  wherein n is 0;
  - Y is selected from the group consisting of  $CR^1$  and  $C(R^2)(R^3)$  and q is 2 or 3.
- 9. A compound of claim 7 which is a derivative thereof selected from the group30 consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

## 10. A compound of claim 7 wherein

is selected from the group consisting of

$$\mathbb{R}^{18} \xrightarrow{\mathbb{N}} \mathbb{N}$$

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wherein R<sup>19</sup>, R<sup>20</sup>, R<sup>21</sup> and R<sup>28</sup> at each occurrence are independently selected from the group consisting of halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -OH, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)(O)NH(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl,

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sulfonyl,  $-SO_2$ - $(C_1-C_3$  alkyl),  $-SO_3$ - $(C_1-C_3$  alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups;

R<sup>18</sup> is selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

R<sup>22</sup> is selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub>alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and

c is an integer of zero to two; d is an integer of zero to three; e is an integer of zero to four; and i is an integer of zero to two.

11. The compound of claim 7 wherein R<sup>18</sup> is aralkyl;

-C(O)NH(benzyl) groups;

 $R^4$  is aryl; T is  $(CH_2)_b$  where b is zero;

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L is (CH<sub>2</sub>)<sub>n</sub> where n is zero; and, B, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup> and R<sup>10</sup> are each independently hydrogen.

## 12. A compound of the structure

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wherein T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of from 0 to 3;

B, R<sup>4</sup>, R<sup>9</sup>, R<sup>10</sup> and R<sup>23</sup> at each occurrence are independently selected from the

L is selected from the group consisting of O, NR<sup>11</sup>, S, and (CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

g is an integer of from 0 to 7; and

group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl,

-CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino,
alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub>
alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)C(O)NH(C<sub>1</sub>-C<sub>3</sub>alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub>
alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino,
alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl,

-C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>,

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-C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and

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-C(O)NH(benzyl) groups;

R<sup>6</sup>, R<sup>7</sup>, R<sup>11</sup> and R<sup>18</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

wherein B, R<sup>4</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>11</sup>, R<sup>18</sup> and R<sup>23</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group;

wherein when L is NR<sup>11</sup>, R<sup>4</sup> and R<sup>11</sup> taken together may form a ring; and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring; or a pharmaceutically acceptable salt thereof.

13. A compound of claim 12 which is a derivative thereof selected from the group

consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

14. A compound of the structure

$$R^{24}$$
 $R^{18}$ 
 $R^{18}$ 
 $R^{18}$ 
 $R^{18}$ 
 $R^{18}$ 
 $R^{18}$ 
 $R^{10}$ 
 $R$ 

wherein h is an integer of zero to five;

B, R<sup>9</sup>, R<sup>10</sup>, R<sup>24</sup> and R<sup>25</sup> are each independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl,
-CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, -OH, alkynylamino, alkoxycarbonyl,

heterocycloyl, carboxy,  $-N(C_1-C_3 \text{ alkyl})-C(O)(C_1-C_3 \text{ alkyl})$ ,

 $-NHC(O)N(C_1-C_3 \text{ alkyl})C(O)NH(C_1-C_3 \text{ alkyl}), -NHC(O)NH(C_1-C_6 \text{ alkyl}),$ -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-5  $C_3$ )alkyl,  $-C(O)N(C_1-C_3 \text{ alkyl})_2$ , -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-10 C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; R<sup>27</sup>, at each occurrence, is independently selected from the group consisting of halogen, hydroxyl, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, 15 -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy,  $-N(C_1-C_3 \text{ alkyl})-C(O)(C_1-C_3 \text{ alkyl})$ ,  $-NHC(O)N(C_1-C_3 \text{ alkyl})C(O)NH(C_1-C_3 \text{ alkyl}), -NHC(O)NH(C_1-C_6 \text{ alkyl}),$ -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), -N(C<sub>1</sub>-C<sub>3</sub>alkyl)SO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub>alkyl), -N(C<sub>1</sub>-C<sub>3</sub>alkyl)SO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-20  $C_3$ )amino,  $-C(O)O-(C_1-C_3)$ alkyl,  $-C(O)NH-(C_1-C_3)alkyl, -C(O)N(C_1-C_3)alkyl, -CH=NOH, -PO_3H_2,$ -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, 25 -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and –C(O)NH(benzyl) groups; R<sup>6</sup>, R<sup>7</sup> and R<sup>18</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, 30 alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy,

carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl,

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cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups; and, R<sup>26</sup> is selected from the group consisting of hydrogen, alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, alkoxycarbonyl, heterocycloyl, carboxy,  $-C(O)O-(C_1-C_3)$  alkyl,  $-C(O)NH-(C_1-C_3)$  alkyl,  $-C(O)N(C_1-C_3)$  alkyl),  $-C(O)N(C_1-C_3)$ PO<sub>3</sub>H<sub>2</sub>, haloalkyl, carboxamide, cycloalkyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, biaryl, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein B,  $R^6$ ,  $R^7$ ,  $R^9$ ,  $R^{10}$ ,  $R^{18}$ ,  $R^{24}$ ,  $R^{25}$ ,  $R^{26}$  and  $R^{27}$  are unsubstituted or substituted with at least one electron donating or electron withdrawing group; wherein R<sup>18</sup> and R<sup>24</sup> taken together may form a ring; R<sup>24</sup> and R<sup>25</sup> taken together may form a ring; R<sup>25</sup> and R<sup>26</sup> taken together may form a ring; and wherein R<sup>9</sup> and R<sup>10</sup> taken together may form a ring: or a pharmaceutically acceptable salt thereof.

- 15. The compound of claim 14 wherein B, R<sup>6</sup>, R<sup>7</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>24</sup>, R<sup>25</sup> and R<sup>26</sup> are each independently hydrogen and R<sup>18</sup> is substituted or unsubstituted aralkyl.
- 16. A compound of claim 14 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.

## 17. A compound of the structure

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**15** 

wherein Z, at each occurrence, is independently selected from the group consisting of C(O), N, CR<sup>30</sup>, C(R<sup>31</sup>)(R<sup>32</sup>), NR<sup>33</sup>, CH, O and S;

z is an integer of from 3 to 6;

k is an integer of from 0 to 5;

T is selected from the group consisting of C(O) and (CH<sub>2</sub>)<sub>b</sub> wherein b is an integer of from 0 to 3;

L is selected from the group consisting of O, NR<sup>11</sup>, S, and

(CH<sub>2</sub>)<sub>n</sub> wherein n is an integer of 0 or 1;

R<sup>6</sup>, R<sup>7</sup>, R<sup>11</sup>, R<sup>18</sup> and R<sup>33</sup> are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

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B, R<sup>4</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>30</sup>, R<sup>31</sup> and R<sup>32</sup> at each occurrence are independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -OH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl),

or a pharmaceutically acceptable salt thereof.

	$-NHC(O)N(C_1-C_3 \ alkyl)C(O)NH(C_1-C_3 alkyl), \ -NHC(O)NH(C_1-C_6 \ alkyl), \\$
	-NHSO <sub>2</sub> (C <sub>1</sub> -C <sub>3</sub> alkyl), -NHSO <sub>2</sub> (aryl), alkoxyalkyl, alkylamino,
	alkenylamino, di(C <sub>1</sub> -C <sub>3</sub> )amino, -C(O)O-(C <sub>1</sub> -C <sub>3</sub> )alkyl, -C(O)NH-(C <sub>1</sub> -
	$C_3$ ) alkyl, $-C(O)N(C_1-C_3$ alkyl) <sub>2</sub> , $-CH=NOH$ , $-PO_3H_2$ , $-OPO_3H_2$ , haloalkyl,
	alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl,
	cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl,
	thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl,
	alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO <sub>2</sub> -(C <sub>1</sub> -C <sub>3</sub> alkyl), -SO <sub>3</sub> -
	(C1-C3 alkyl), sulfonamido, carbamate, aryloxyalkyl and
	-C(O)NH(benzyl) groups; and
R <sup>29</sup> , at o	each occurrence, is independently selected from the group consisting of
	halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy,
	hydroxyalkyl, aliphatic acyl, -CF <sub>3</sub> , -CO <sub>2</sub> H, -SH, -CN, -NO <sub>2</sub> , -NH <sub>2</sub> , -OH,
	alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C <sub>1</sub> -C <sub>3</sub> alkyl)-
	$C(O)(C_1-C_3 \text{ alkyl}), -NHC(O)N(C_1-C_3 \text{ alkyl})C(O)NH(C_1-C_3 \text{alkyl}),$
	-NHC(O)NH(C <sub>1</sub> -C <sub>6</sub> alkyl), -NHSO <sub>2</sub> (C <sub>1</sub> -C <sub>3</sub> alkyl), -NHSO <sub>2</sub> (aryl),
	alkoxyalkyl, alkylamino, alkenylamino, di(C1-C3)amino, -C(O)O-(C1-
	$C_3$ )alkyl, $-C(O)NH-(C_1-C_3)$ alkyl, $-C(O)N(C_1-C_3)$ alkyl) <sub>2</sub> , $-CH=NOH$ ,
	$-PO_3H_2,-OPO_3H_2,haloalkyl,alkoxyalkoxy,carboxaldehyde,carboxamide$
	cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl,
	aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl,
	aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl,
	-SO <sub>2</sub> -(C <sub>1</sub> -C <sub>3</sub> alkyl), -SO <sub>3</sub> -(C <sub>1</sub> -C <sub>3</sub> alkyl), sulfonamido, carbamate,
	aryloxyalkyl and -C(O)NH(benzyl) groups;
wherein	n B, $R^4$ , $R^5$ , $R^6$ , $R^7$ , $R^9$ , $R^{10}$ , $R^{11}$ , $R^{18}$ , $R^{29}$ , $R^{30}$ , $R^{31}$ , $R^{32}$ and $R^{33}$ are
	unsubstituted or substituted with at least one electron donating or electron
	withdrawing group;
	wherein when L is NR <sup>11</sup> , R <sup>4</sup> and R <sup>11</sup> taken together may form a ring;
	and wherein R <sup>9</sup> and R <sup>10</sup> taken together may form a ring;

- 18. A compound of claim 17 which is a derivative thereof selected from the group consisting of esters, carbamates, aminals, amides, optical isomers and pro-drugs.
- The compound of claim 17 wherein z is three or four. 19.

20. A compound of the structure

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wherein  $R^{24}$  and  $R^{25}$  are each independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF3, -SH, -OH, -CO<sub>2</sub>H, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy,  $-N(C_1-C_3 \text{ alkyl})-C(O)(C_1-C_3 \text{ alkyl})$ ,  $-NHC(O)N(C_1-C_3 \text{ alkyl})$ alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>- $C_3$ )amino, -C(O)O-( $C_1$ - $C_3$ )alkyl, -C(O)NH-( $C_1$ - $C_3$ )alkyl, - $C(O)N(C_1$ - $C_3$ alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; and

 $R^{18}$  and  $R^{34}$  are each independently selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde,

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carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups;

wherein R<sup>18</sup>, R<sup>24</sup>, R<sup>25</sup> and R<sup>34</sup> are unsubstituted or substituted with at least one electron donating or electron withdrawing group;

and wherein  $R^{24}$  and  $R^{25}$  taken together may form a ring; with the proviso that when  $R^{24}$  and  $R^{25}$  taken together form a ring, the ring formed is not benzene.

21. A compound of claim 20 wherein R<sup>34</sup> is hydrogen;

 $R^{18}$  is aralkyl; and  $R^{24}$  and  $R^{25}$  are each independently selected from the group consisting of hydrogen, lower alkyl, and lower alkyl wherein  $R^{24}$  and  $R^{25}$  taken together may form a ring.

22. A compound of claim 20 of the structure

$$\mathbb{R}^{24}$$
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 
 $\mathbb{R}^{24}$ 

wherein R<sup>24</sup> and R<sup>25</sup> are each independently selected from the group consisting of hydrogen, halogen, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -SH, -OH, -CO<sub>2</sub>H, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy, -N(C<sub>1</sub>-C<sub>3</sub> alkyl)-C(O)(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHC(O)NH(C<sub>1</sub>-C<sub>6</sub> alkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub>

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alkyl), -NHSO2(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C1-

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C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)N(C<sub>1</sub>-C<sub>3</sub> alkyl)2, -CH=NOH, -PO3H2, -OPO3H2, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; R<sup>34</sup> is selected from the group consisting of alkyl, alkenyl, alkynyl, hydroxyalkyl, aliphatic acyl, alkynylamino, alkoxycarbonyl, heterocycloyl, -CH=NOH, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, carbamate, aryloxyalkyl, hydrogen and -C(O)NH(benzyl) groups; and, R<sup>35</sup>, at each occurrence, is independently selected from the group consisting of halogen, hydroxyl, alkyl, alkenyl, alkynyl, alkoxy, alkenoxy, alkynoxy, thioalkoxy, hydroxyalkyl, aliphatic acyl, -CF<sub>3</sub>, -CO<sub>2</sub>H, -SH, -CN, -NO<sub>2</sub>, -NH<sub>2</sub>, alkynylamino, alkoxycarbonyl, heterocycloyl, carboxy,  $-N(C_1-C_3 \text{ alkyl})-C(O)(C_1-C_3 \text{ alkyl})$ ,  $-NHC(O)N(C_1-C_3 \ alkyl)C(O)NH(C_1-C_3 alkyl), \ -NHC(O)NH(C_1-C_6 \ alkyl), \\$ -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>3</sub> alkyl), -NHSO<sub>2</sub>(aryl), alkoxyalkyl, alkylamino, alkenylamino, di(C<sub>1</sub>-C<sub>3</sub>)amino, -C(O)O-(C<sub>1</sub>-C<sub>3</sub>)alkyl, -C(O)NH-( $C_1$ - $C_3$ )alkyl, -C(O)N( $C_1$ - $C_3$  alkyl)<sub>2</sub>, -CH=NOH, -PO<sub>3</sub>H<sub>2</sub>, -OPO<sub>3</sub>H<sub>2</sub>, haloalkyl, alkoxyalkoxy, carboxaldehyde, carboxamide, cycloalkyl, cycloalkenyl, cycloalkynyl, cycloalkylalkyl, aryl, aroyl, aryloxy, arylamino, biaryl, thioaryl, diarylamino, heterocyclyl, alkylaryl, aralkenyl, aralkyl, alkylheterocyclyl, heterocyclylalkyl, sulfonyl, -SO<sub>2</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), -SO<sub>3</sub>-(C<sub>1</sub>-C<sub>3</sub> alkyl), sulfonamido, carbamate, aryloxyalkyl and -C(O)NH(benzyl) groups; wherein  $R^{24}$ ,  $R^{25}$ ,  $R^{34}$  and  $R^{35}$  are unsubstituted or substituted with

at least one electron donating or electron withdrawing group; and,

m is an integer of from 0 to 5.

5 23. A compound of claim 22 wherein R<sup>34</sup> is hydrogen;

m is an integer of one to three and R<sup>35</sup> at each occurrence is selected from the group consisting of alkyl, halogen, alkoxy, haloalkyl, sulfonyl, -OH and -CN.

- 24. A compound of claim 20 selected from the group consisting of
- 5-(2-chlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-fluorobenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-fluorobenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-benzyl-6-methyl-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-benzyl-3,5-dihydro[1,3]oxazolo[4,5-
  - c|pyridine-2,4-dione, 5-(2,5-dimethylbenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4
    - dione, 5-(2-methylbenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,4-
    - dichlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-methoxybenzyl)-3,5-
    - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,5-difluorobenzyl)-3,5-
    - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(methylthio)benzyl]-3,5-
- 20 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-fluorobenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-methoxybenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3,5-bis(trifluoromethyl)benzyl]-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-tert-butylbenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-chlorobenzyl)-3,5-
- 25 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chlorobenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-(trifluoromethyl)benzyl]-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-bromobenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3,4-dichlorobenzyl)-3,5-
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-methylbenzyl)-3,5-
- 30 dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methoxybenzyl)-3,5
  - dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[4-(trifluoromethyl)benzyl]-3,5-

- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-methylbenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(pyridin-2-ylmethyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,4-difluorobenzyl)-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-difluorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-(trifluoromethoxy)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[4-(trifluoromethoxy)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(trifluoromethyl)benzyl]-3,5-

dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-methoxybenzyl)-3,5-

- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,3-dichlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3,5-dimethylbenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-pentyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,4-dichlorobenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-ethyl-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-butyl-5-(2-chlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(trifluoromethyl)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-dichlorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-fluorobenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methylbenzyl)-7-methyl-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chlorobenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 7-methyl-5-[4-(methylsulfonyl)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-methoxybenzyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-
- propyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 4-[(2,4-dioxo-2,3-dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]-N,N-dimethylbenzenesulfonamide, 5-(mesitylmethyl)-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-3,5,6,7,8,9-hexahydro[1,3]oxazolo[4,5-c]quinoline-2,4-dione, 5-(2-chlorobenzyl)-7-ethyl-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(methylthio)benzyl]-3,5-
- dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 2-[(2,4-dioxo-2,3-dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]-N,N-dimethylbenzenesulfonamide, 5-(2,6-dimethoxybenzyl)-3,5-

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dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(trifluoromethoxy)benzyl]-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-6,7-dimethyl-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(methylsulfonyl)benzyl]-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chloro-2-methoxybenzyl)-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-5,6,7,8,9,10-hexahydro-
     2H-cyclohepta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-[2-(difluoromethoxy)benzyl]-
     3.5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-methyl-5-[(1R)-1-phenylethyl]-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(4-chlorobenzyl)-7-propyl-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(methylsulfonyl)benzyl]-3,5-
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     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-dimethylbenzyl)-3,5-
     dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 3-chloro-2-[(2,4-dioxo-2,3-
      dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]benzonitrile, 5-(2-chloro-6-
      methylbenzyl)-6,7-dimethyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 2-[(2,4-dioxo-
      2.3-dihydro[1,3]oxazolo[4,5-c]pyridin-5(4H)-yl)methyl]benzonitrile, 5-(2-chloro-6-
      methoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-
15
      (methylthio)benzyl]-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-
      cyclopropyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3-chlorobenzyl)-7-methyl-
      3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-dichlorobenzyl)-7-methyl-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-methyl-5-(4-methylbenzyl)-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(3,5-dimethoxybenzyl)-7-methyl-3,5-
20
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,6-difluorobenzyl)-7-methyl-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[3-(methylsulfonyl)benzyl]-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-7-methyl-3,5-
25
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-fluoro-6-methoxybenzył)-7-methyl-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methoxybenzyl)-7-propyl-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(5-chloro-2-fluorobenzyl)-7-methyl-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-isopropyl-3,5-
      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(5-fluoro-2-methylbenzyl)-7-methyl-3,5-
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      dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 7-methyl-5-[(1S)-1-phenylethyl]-3,5-
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dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-isopropoxybenzyl)-7-methyl-

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3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(5-acetyl-2-methoxybenzyl)-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chlorobenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-d]pyridazine-2,4-dione, 5-[2-fluoro-6-(trifluoromethyl)benzyl]-7methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-methylbenzyl)-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-(2-chloro-6-ethoxybenzyl)-7-ethyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6propoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6isobutoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)dione, 5-(2-chloro-6-isopropoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-6-(2,2,2-trifluoroethoxy)benzyl]-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-7-methyl-3,5dihydro[1,3]oxazolo[4,5-d]pyridazine-2,4-dione, 5-[2-chloro-6-(2-methoxyethoxy)benzyl]-5,6,7,8-tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-(2-chloro-6-ethoxybenzyl)-6,7-dimethyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6-ethoxybenzyl)-7-ethyl-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2chlorobenzyl)-7-ethyl-3,5-dihydro[1,3]oxazolo[4,5-d]pyridazine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-7-propyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-7-cyclopropyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5-propoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5methoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-6ethoxybenzyl)-6-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2-chloro-5ethoxybenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(piperidin-1-ylsulfonyl)benzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-chloro-5-(pyrrolidin-1-ylsulfonyl)benzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5c]pyridine-2,4-dione, 5-[2-chloro-6-(cyclopentylmethoxy)benzyl]-7-methyl-3,5dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-[2-(benzyloxy)-6-chlorobenzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione, 5-(2,3-dichloro-6-ethoxybenzyl)-5,6,7,8tetrahydro-2H-cyclopenta[b][1,3]oxazolo[5,4-d]pyridine-2,4(3H)-dione, 5-[2-chloro-5-(trifluoromethyl)benzyl]-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione and 5-

(2-chloro-5-fluorobenzyl)-7-methyl-3,5-dihydro[1,3]oxazolo[4,5-c]pyridine-2,4-dione.

- 25. A compound selected from the group consisting of (3S)-3-[({[2-methyl-4-(2-methylpropyl)-6-oxo-1-(phenylmethyl)-1,6-dihydro-5-pyrimidinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-(1,3-dihydro-5-pyrimidinyl)amino]-3-(4-methylphenylphenyl
- benzodioxol-5-yl)-3-[({[2-oxo-1-(phenylmethyl)-4-propyl-1,2-dihydro-3-pyridinyl]amino}carbonyl)amino]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-ethyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-2-oxo-4-propyl-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid,
- 10 (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({6-methyl-2-oxo-1-(phenylmethyl)-4-[(phenylmethyl)oxy]-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-2,4-dimethyl-6-oxo-1,6-dihydro-5-yrimidinyl}amino)carbonyl]amino}-
- 3-(4-methylphenyl)propanoic acid, (3S)-3-{[({4-amino-1-[(2-chlorophenyl)methyl]-6-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[4-(methyloxy)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-
- 20 (3,4-dimethylphenyl)propanoic acid, (3S)-3-{[({4-amino-1-[(2-chlorophenyl)methyl]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-[(2-chlorophenyl)methyl]-4-(1,4-oxazinan-4-yl)-2-oxo-1,2-dihydro-3-
- pyridinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-[(2-chlorophenyl)methyl]-2-oxo-4-(propylamino)-1,2-dihydro-3-pyridinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-bromophenyl)methyl]-4-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}
- dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[3-methyl-4-(methyloxy)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-2-oxo-4-phenyl-1,2-dihydro-3-

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pyridinyl\} amino) carbonyl] amino\} -3-(4-methylphenyl) propanoic acid, (3S)-3-\{[(\{1-[(2-chlorophenyl)methyl]-4-[(2-\{[2-(methyloxy)ethyl]oxy\}ethyl)oxy]-2-oxo-1,2-dihydro-3-pyridinyl\} amino) carbonyl] amino\} -3-(4-methylphenyl) propanoic acid, (3S)-3-\{[(\{1-[(2-chlorophenyl)methyl]-4-hydroxy-6-methyl-2-oxo-1,2-dihydro-3-
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- pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-[(1,1-dimethylethyl)amino]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-phenylpropanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-[4-methyltetrahydro-1(2H)-4-[4-methyltetrahydro-1(4-methylte
- pyrazinyl]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[4-(methyloxy)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-
- pyridinyl}amino)carbonyl]amino}-3-(3,5-dimethylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(3-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[3-(methyloxy)phenyl]propanoic acid, (3S)-3-[3,5-bis(methyloxy)phenyl]-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}
- chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-quinolinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-[3-(trifluoromethyl)phenyl]propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-({ethyl[(ethylamino)carbonyl]amino}carbonyl)amino]-2-oxo-1,2-dihydro-3-
- pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({4-(1-azetanyl)-1-[(2-chlorophenyl)methyl]-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-[(2-chlorophenyl)methyl]-4-({2-[(2-{[2-(methyloxy)ethyl]oxy}ethyl)oxy]ethyl}oxy)-2-oxo-1,2-dihydro-3-pyridinyl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-
- 30 {[({1-[(2-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl}amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2-fluorophenyl)methylphenyl]-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl}amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-(4-methylphenyl)propanoic acid, (3S)-3-(4-methylphenylp

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chloro-6-fluorophenyl)methyl]-4-hydroxy-2-oxo-1,2-dihydro-3pyridinyl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[(2chlorophenyl)methyl]-5-methyl-2-oxo-1,2-dihydro-3-pyridinyl}amino)carbonyl]amino}-3-(4-(trifluoromethyl)phenyl)methyl)-1,2 dihydro-3-pyridinyl)amino)carbonyl)amino)propanoic acid, (3S)-3-((((1-((2-chlorophenyl)methyl)-2-oxo-1,2-dihydro-3-fluorophenyl)methyl)-2-oxo-1,2-dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4methylphenyl)propanoic acid, (3S)-3-((((1-((2-bromophenyl)methyl)-2-oxo-1,2-dihydro-3pyridinyl)amino)carbonyl)amino)-3-(4-methylphenyl)propanoic acid, (3S)-3-((((1-((2,4dichlorophenyl)methyl)-2-oxo-1,2-dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4methylphenyl)propanoic acid, (3S)-3-((((1-((2-chloro-6-fluorophenyl)methyl)-2-oxo-1,2dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4-methylphenyl)propanoic acid, (3S)-3-((((1-((2-chlorophenyl)methyl)-4-hydroxy-2-oxo-1,2-dihydro-3-pyridinyl)amino)carbonyl)amino)-3-(4-trifluoromethyl)oxy)phenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-methoxybenzyl)-2oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, 4-{[3-[({[(1S)-2-carboxy-1-(4-methylphenyl)ethyl]amino}carbonyl)amino]-1-(2-chlorobenzyl)-2-oxo-1,2-dihydropyridin-4-yl]amino}benzoic acid, (3S)-3-{[({1-(2-chlorobenzyl)-4-[(2,2dimethylpropanoyl)amino]-2-oxo-1,2-dihydropyridin-3-yl}amino)carbonyl]amino}-3-(4methylphenyl)propanoic acid, (3S)-3-[({[4-{[(tert-butylamino)carbonyl]amino}-1-(2chlorobenzyl)-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4methylphenyl)propanoic acid, (3S)-3-[({[1-(2-cyanobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(2,3dihydro-1,4-benzodioxin-6-yl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(7-methoxy-1,3-benzodioxol-5yl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(3-ethoxy-4-methoxyphenyl)propanoic acid, (3S)-3-[({[1-(2chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4dimethoxyphenyl)propanoic acid, (3S)-3-[({[1-(4-chlorobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-

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(2-chloro-6-methoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
          yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-
          4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-
          methylphenyl)propanoic acid, (3S)-3-[({[1-(2,6-difluorobenzyl)-4-hydroxy-2-oxo-1,2-
          dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-
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          (2-chloro-6-methoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
          yl]amino}carbonyl)amino]-3-(3,5-dimethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-
          chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino]carbonyl)amino]-3-(3,4-
          diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-
          dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-
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          (2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-
          methoxy-4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-
          1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,5-dimethoxy-4-
          methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-
          dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-dimethylphenyl)propanoic acid, (3S)-3-
15
          [({[1-(2-chlorobenzyl)-5-ethyl-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
          yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-{[({1-[2-chloro-5-
           (trifluoromethyl)benzyl]-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl}amino)carbonyl]amino}-
           3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-methoxybenzyl)-4-hydroxy-2-
           oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-methylphenyl)propanoic acid,
20
           (3S)-3-[(\{[1-(2-chloro-6-methylbenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-methyl-2-oxo-1,2-dihydropyridin-3-methylbenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-methyl-2-oxo-1,2-dihydropyridin-3-methylbenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-methyl-2-oxo-1,2-dihydropyridin-3-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihydroxy-5-methyl-2-oxo-1,2-dihy
           yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-
           4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-
           (4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2,6-dimethoxybenzyl)-4-hydroxy-2-oxo-1,2-
           dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-
25
           (2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-
           propoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-5-propyl-1,2-
           dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-
           (2-chlorobenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-
           yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-
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4-hydroxy-2-oxo-5-propyl-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-

diethoxyphenyl)propanoic acid, (3S)-3-(3-butoxyphenyl)-3-[({[1-(2-chlorobenzyl)-4hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]propanoic acid, (3S)-3-{[({1-[2-chloro-5-(methylsulfonyl)benzyl]-4-hydroxy-2-oxo-1,2-dihydropyridin-3yl}amino)carbonyl]amino}-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(2-5 methoxyethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-dipropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(difluoromethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-10 diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-methylbenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(3,4-diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6cyanobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-15 methylphenyl)propanoic acid, 3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(2-naphthyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4diethoxyphenyl)propanoic acid (3S)-3-[({[1-(2-chloro-6-methoxybenzyl)-4-hydroxy-5methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3,4-20 diethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-1,2dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3yl]amino}carbonyl)amino]-3-(4-methoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6methylbenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-25 yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6ethoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-30 cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-

- [({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(1-methyl-1H-indol-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(2,3-dihydro-1-benzofuran-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-
- hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3,5-diethoxyphenyl)propanoic acid, (3S)-3-[({[5-chloro-1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-
- [({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-propoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-phenylpropanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-
- yl]amino}carbonyl)amino]-3-(1,3-diethyl-2-oxo-2,3-dihydro-1H-benzimidazol-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(trifluoromethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5,6-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dimethyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-dimethyl-2-oxo-1,2-dihydropyridin-3-dihydropyridin-3-dihydroyy-1-dihyd
- chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(1-methyl-1H-indol-5-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-5-cyclopropyl-4-hydroxy-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-5-cyclopropyl-4-hydroxy-2-oxo-1,2-dihydropyridin-3-
- yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-5-methoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-6-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-eth
- methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(1-methyl-1H-indol-6-yl)propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-2-oxo-2,5,6,7-

tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-[3-(cyclopropyloxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-[3-(cyclopropylmethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-

- hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-[3-(cyclopropylmethoxy)phenyl]propanoic acid, (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3,5-dimethylphenyl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-
- [(difluoromethyl)oxy]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-[(1,1,2,2-tetrafluoroethyl)oxy]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-(1-ethyl-1H-indol-5-yl)propanoic acid and (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-[3-(diethylamino)phenyl]propanoic acid and pharmaceutical acceptable salts thereof.
  - 26. (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid and pharmaceutically acceptable salts thereof.
    - 27. (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid and pharmaceutically acceptable salts thereof.
    - 28. (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-[3-(diethylamino)phenyl]propanoic acid and pharmaceutically acceptable salts thereof.

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- 29. A compound selected from the group consisting of (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(4-methylphenyl)propanoic acid; (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-
- methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(3-ethoxyphenyl)propanoic acid; (3S)-3-[({[1-(2-chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-isopropoxyphenyl)propanoic acid; (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(6-methoxy-2-naphthyl)propanoic acid; (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl-2-oxo-1,2-dihydropyridin-3-yl]amino}carbonyl)amino]-3-(6-methoxy-2-naphthyl)propanoic acid; (3S)-3-[({[1-(2-chloro-6-ethoxybenzyl)-4-hydroxy-5-methyl)propanoic acid; (3S)-3-[({[1-(2-chl
- chlorobenzyl)-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl]amino}carbonyl)amino]-3-(3-methylphenyl)propanoic acid; (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-(1-methyl-1H-indol-5-yl)propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-
- yl}amino)carbonyl]amino}-3-{3-[(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{[(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-
- 20 {3-[methyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-[methyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chlorophenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-
- 25 [ethyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-{3-[ethyl(methylsulfonyl)amino]phenyl}propanoic acid, (3S)-3-{[({1-[(2-chloro-6-methylphenyl)methyl]-4-hydroxy-2-oxo-2,5,6,7-tetrahydro-1H-cyclopenta[b]pyridin-3-yl}amino)carbonyl]amino}-3-(1H-indol-5-yl)propanoic acid and pharmaceutically acceptable salts thereof.

- 30. A pharmaceutical composition comprising:a compound of claim 1in a pharmaceutically acceptable carrier.
- 5 31. A method for selectively inhibiting  $\alpha_4\beta_1$  integrin binding in a mammal comprising administering to said mammal a therapeutic amount of a compound of claim 1.